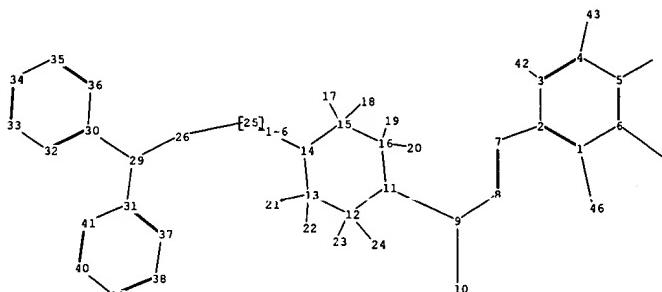
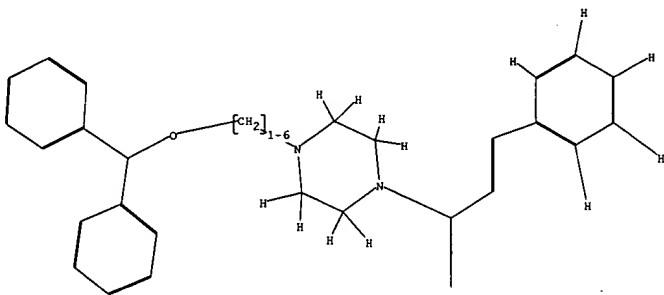


update  
4/10/53



chain nodes :

7 8 9 17 18 19 20 21 22 23 24 25 26 29 42 43 44 45 46

ring nodes :

1 2 3 4 5 6 11 12 13 14 15 16 30 31 32 33 34 35 36 37 38 39 40 41

ring/chain nodes :

10

chain bonds :

1-46 2-7 3-42 4-43 5-44 6-45 7-8 8-9 9-10 9-11 12-23 12-24 13-21 13-22  
14-25 15-17 15-18 16-19 16-20 25-26 26-29 29-30 29-31

ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6 11-12 11-16 12-13 13-14 14-15 15-16 30-32 30-36  
31-37 31-41 32-33 33-34 34-35 35-36 37-38 38-39 39-40 40-41

exact/norm bonds :

9-11 11-12 11-16 12-13 13-14 14-15 15-16 26-29

exact bonds :

1-46 2-7 3-42 4-43 5-44 6-45 7-8 8-9 9-10 12-23 12-24 13-21 13-22 14-25  
15-17 15-18 16-19 16-20 25-26 29-30 29-31

normalized bonds :

1-2 1-6 2-3 3-4 4-5 5-6 30-32 30-36 31-37 31-41 32-33 33-34 34-35 35-36  
37-38 38-39 39-40 40-41

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:CLASS 8:CLASS 9:CLASS 10:CLASS  
11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:CLASS 18:CLASS 19:CLASS  
20:CLASS 21:CLASS 22:CLASS 23:CLASS 24:CLASS 25:CLASS 26:CLASS 29:CLASS 30:Atom  
31:Atom 32:Atom 33:Atom 34:Atom 35:Atom 36:Atom 37:Atom 38:Atom 39:Atom 40:Atom  
41:Atom 42:CLASS 43:CLASS 44:CLASS 45:CLASS 46:CLASS

=> S L1 FULL  
 FULL SEARCH INITIATED 22:19:48 FILE 'REGISTRY'  
 FULL SCREEN SEARCH COMPLETED - 9 TO ITERATE

100.0% PROCESSED 9 ITERATIONS  
 SEARCH TIME: 00.00.01

2 ANSWERS

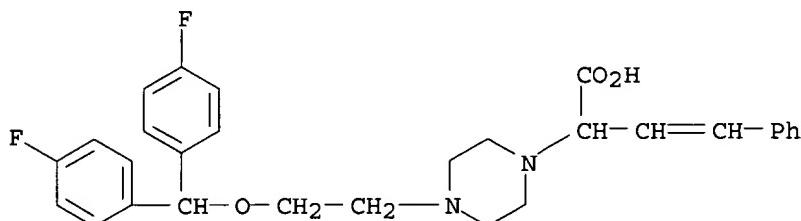
L3 2 SEA SSS FUL L1

=> D ALL

L3 ANSWER 1 OF 2 REGISTRY COPYRIGHT 2003 ACS on STN  
 RN 350583-56-1 REGISTRY  
 CN 1-Piperazineacetic acid,  
 4-[2-[bis(4-fluorophenyl)methoxy]ethyl]-.alpha.-(2-phenylethenyl)- (9CI) (CA  
 INDEX NAME)  
 FS 3D CONCORD  
 MF C29 H30 F2 N2 O3  
 SR CA  
 LC STN Files: CA, CAPLUS, USPATFULL

#### Ring System Data

Elemental Analysis	Elemental Sequence	Size of the Rings	Ring System Formula	Ring Identifier RID	RID Occurrence Count
EA	ES	SZ	RF	RID	Count
C6	C6	6	C6	46.150.18	3
C4N2	NC2NC2	6	C4N2	46.383.1	1



#### Calculated Properties (CALC)

PROPERTY (CODE)	VALUE	CONDITION	NOTE
Bioconc. Factor (BCF)	18.0	pH 1	(1) ACD
Bioconc. Factor (BCF)	137	pH 4	(1) ACD
Bioconc. Factor (BCF)	160	pH 7	(1) ACD
Bioconc. Factor (BCF)	125	pH 8	(1) ACD
Bioconc. Factor (BCF)	8.85	pH 10	(1) ACD
Boiling Point (BP)	623.1+/-55.0 deg C	760.0 Torr	(1) ACD
Enthalpy of Vap. (HVAP)	97.02+/-3.0 kJ/mol		(1) ACD
Flash Point (FP)	330.6+/-56.7 deg C		(1) ACD
H acceptors (HAC)	5		(1) ACD
H donors (HD)	1		(1) ACD
Koc (KOC)	28.6	pH 1	(1) ACD
Koc (KOC)	218	pH 4	(1) ACD
Koc (KOC)	255	pH 7	(1) ACD
Koc (KOC)	200	pH 8	(1) ACD

Koc (KOC)	14.1	pH 10	(1) ACD
logD (LOGD)	3.05	pH 1	(1) ACD
logD (LOGD)	3.93	pH 4	(1) ACD
logD (LOGD)	4.00	pH 7	(1) ACD
logD (LOGD)	3.89	pH 8	(1) ACD
logD (LOGD)	2.74	pH 10	(1) ACD
logP (LOGP)	6.503+/-0.668		(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 1	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 4	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 7	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 8	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 10	(1) ACD
Molecular Weight (MW)	492.56		(1) ACD
pKa (PKA)	3.93+/-0.10	Most Acidic	(1) ACD
pKa (PKA)	8.47+/-0.50	Most Basic	(1) ACD
Vapor Pressure (VP)	2.19E-16 Torr	25.0 deg C	(1) ACD

(1) Calculated using Advanced Chemistry Development (ACD) Software Solaris V4.67  
((C) 1994-2003 ACD)

1 REFERENCES IN FILE CA (1907 TO DATE)  
1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

#### REFERENCE 1

AN 135:107255 CA  
TI Preparation of polypharmacophoric agents  
IN Hanson, Robert N.; Babich, John W.  
PA Biostream Therapeutics, Inc., USA  
SO PCT Int. Appl., 74 pp.  
CODEN: PIXXD2  
DT Patent  
LA English  
IC ICM C07D241-04  
ICS C07D211-34; C07D211-44; C07D223-26  
CC 27-16 (Heterocyclic Compounds (One Hetero Atom))  
Section cross-reference(s): 1  
FAN.CNT 1  
PATENT NO. KIND DATE APPLICATION NO. DATE  
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PI WO 2001051474 A2 20010719 WO 2001-US1035 20010111  
WO 2001051474 A3 20011206  
W: CA, JP  
RW: AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT,  
SE, TR  
US 2002042357 A1 20020411 US 2001-758957 20010111  
EP 1257541 A2 20021120 EP 2001-902026 20010111  
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE,  
FI, CY, TR  
JP 2003519689 T2 20030624 JP 2001-551856 20010111  
PRAI US 2000-175617P 20000111  
WO 2001-US1035 20010111  
AB Title compds. (I) comprise a scaffold bearing .gtoreq.2 pharmacophore units selected from D1, D2, D3, and D4 agonists, (ir)reversible monoamine inhibitors, monoamine transporter inhibitors, COMT inhibitors, MAO inhibitors, and dopamine transporter inhibitors. I interact with .gtoreq.2 biol. targets. Thus, (E)-PhZCH(CO<sub>2</sub>H)CH:CHPh (Z = piperidine-4,1-diyl) was prep'd. Data for biol. activity of I were given.  
ST polypharmacophoric agent prepn; dopaminergic system agent prepn  
IT Dopamine agonists  
(D1; prepn. of polypharmacophoric agents)  
IT Dopamine agonists

(D2; prepn. of polypharmacophoric agents)  
 IT Dopamine agonists  
 (D3; prepn. of polypharmacophoric agents)  
 IT Nervous system  
 (Huntington's chorea, treatment; prepn. of polypharmacophoric agents)  
 IT Mental disorder  
 (attention deficit disorder, treatment; prepn. of polypharmacophoric agents)  
 IT Mental disorder  
 (autism, treatment; prepn. of polypharmacophoric agents)  
 IT Transport proteins  
 RL: BPR (Biological process); BSU (Biological study, unclassified); BIOL  
 (Biological study); PROC  
 (Process)  
 (dopamine-transporting, inhibitors; prepn. of polypharmacophoric agents)  
 IT Nervous system  
 (dopaminergic; prepn. of polypharmacophoric agents)  
 IT Monoamines  
 RL: BPR (Biological process); BSU (Biological study, unclassified); BIOL  
 (Biological study); PROC  
 (Process)  
 (inhibitors; prepn. of polypharmacophoric agents)  
 IT Transport proteins  
 RL: BPR (Biological process); BSU (Biological study, unclassified); BIOL  
 (Biological study); PROC  
 (Process)  
 (monoamine-transporting, inhibitors; prepn. of polypharmacophoric agents)  
 IT Anti-inflammatory agents  
 Antidepressants  
 Antiobesity agents  
 Pharmacophores  
 (prepn. of polypharmacophoric agents)  
 IT Alzheimer's disease  
 (treatment; prepn. of polypharmacophoric agents)  
 IT 9001-66-5, monoamine oxidase 9012-25-3, Catechol O-methyl transferase  
 RL: BPR (Biological process); BSU (Biological study, unclassified); BIOL  
 (Biological study); PROC  
 (Process)  
 (inhibitors; prepn. of polypharmacophoric agents)  
 IT 67469-69-6P 350583-53-8P 350583-56-1P 350583-58-3P 350583-59-4P  
 350583-60-7P 350583-61-8P  
 350583-62-9P 350583-63-0P 350583-64-1P 350583-65-2P 350583-66-3P  
 350583-67-4P  
 RL: BAC (Biological activity or effector, except adverse); BSU (Biological  
 study, unclassified); SPN  
 (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP  
 (Preparation); USES (Uses)  
 (prepn. of polypharmacophoric agents)

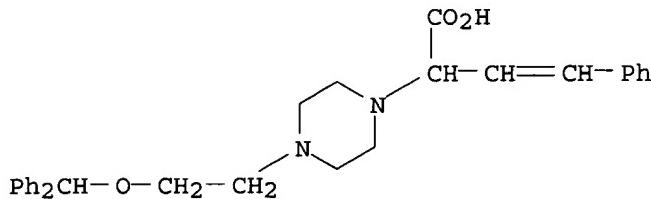
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L3 ANSWER 2 OF 2 REGISTRY COPYRIGHT 2003 ACS on STN  
 RN 350583-53-8 REGISTRY  
 CN 1-Piperazineacetic acid,  
 4-[2-(diphenylmethoxy)ethyl]-.alpha.- (2-phenylethenyl)- (9CI) (CA INDEX NAME)  
 FS 3D CONCORD  
 MF C29 H32 N2 O3  
 SR CA  
 LC STN Files: CA, CAPLUS, USPATFULL

Ring System Data

Elemental Analysis	Elemental Sequence	Size of the Rings	Ring System	Ring Formula	Identifier	RID Occurrence
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EA	ES	SZ	RF	RID	Count
C6	C6	6	C6	46.150.18	3
C4N2	NC2NC2	6	C4N2	46.383.1	1



#### Calculated Properties (CALC)

PROPERTY (CODE)	VALUE	CONDITION	NOTE
Bioconc. Factor (BCF)	14.9	pH 1	(1) ACD
Bioconc. Factor (BCF)	112	pH 4	(1) ACD
Bioconc. Factor (BCF)	134	pH 7	(1) ACD
Bioconc. Factor (BCF)	105	pH 8	(1) ACD
Bioconc. Factor (BCF)	7.50	pH 10	(1) ACD
Boiling Point (BP)	621.5+/-55.0 deg C	760.0 Torr	(1) ACD
Enthalpy of Vap. (HVAP)	96.80+/-3.0 kJ/mol		(1) ACD
Flash Point (FP)	329.6+/-56.7 deg C		(1) ACD
H acceptors (HAC)	5		(1) ACD
H donors (HD)	1		(1) ACD
Koc (KOC)	24.9	pH 1	(1) ACD
Koc (KOC)	188	pH 4	(1) ACD
Koc (KOC)	224	pH 7	(1) ACD
Koc (KOC)	177	pH 8	(1) ACD
Koc (KOC)	12.6	pH 10	(1) ACD
logD (LOGD)	2.94	pH 1	(1) ACD
logD (LOGD)	3.82	pH 4	(1) ACD
logD (LOGD)	3.89	pH 7	(1) ACD
logD (LOGD)	3.79	pH 8	(1) ACD
logD (LOGD)	2.64	pH 10	(1) ACD
logP (LOGP)	6.400+/-0.547		(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 1	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 4	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 7	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 8	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 10	(1) ACD
Molecular Weight (MW)	456.58		(1) ACD
pKa (PKA)	3.93+/-0.10	Most Acidic	(1) ACD
pKa (PKA)	8.48+/-0.50	Most Basic	(1) ACD
Vapor Pressure (VP)	2.64E-16 Torr	25.0 deg C	(1) ACD

(1) Calculated using Advanced Chemistry Development (ACD) Software Solaris V4.67  
 ((C) 1994-2003 ACD)

1 REFERENCES IN FILE CA (1907 TO DATE)  
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REFERENCE 1

AN 135:107255 CA

TI Preparation of polypharmacophoric agents

IN Hanson, Robert N.; Babich, John W.

PA Biostream Therapeutics, Inc., USA

SO PCT Int. Appl., 74 pp.

CODEN: PIXXD2

DT Patent

LA English

IC ICM C07D241-04

ICS C07D211-34; C07D211-44; C07D223-26

CC 27-16 (Heterocyclic Compounds (One Hetero Atom))

Section cross-reference(s): 1

FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2001051474	A2	20010719	WO 2001-US1035	20010111
	WO 2001051474	A3	20011206		
	W: CA, JP				
	RW: AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT,				

SE, TR

US 2002042357	A1	20020411	US 2001-758957	20010111
EP 1257541	A2	20021120	EP 2001-902026	20010111

R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE,

FI, CY, TR

JP 2003519689	T2	20030624	JP 2001-551856	20010111
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PRAI US 2000-175617P 20000111

WO 2001-US1035 20010111

AB Title compds. (I) comprise a scaffold bearing .gtoreq.2 pharmacophore units selected from D1, D2, D3, and D4 agonists, (ir)reversible monoamine inhibitors, monoamine transporter inhibitors, COMT inhibitors,

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(D1; prepn. of polypharmacophoric agents)

IT Dopamine agonists

(D2; prepn. of polypharmacophoric agents)

IT Dopamine agonists

(D3; prepn. of polypharmacophoric agents)

IT Nervous system

(Huntington's chorea, treatment; prepn. of polypharmacophoric agents)

IT Mental disorder

(attention deficit disorder, treatment; prepn. of polypharmacophoric agents)

IT Mental disorder

(autism, treatment; prepn. of polypharmacophoric agents)

IT Transport proteins

RL: BPR (Biological process); BSU (Biological study, unclassified); BIOL (Biological study); PROC (Process)

(dopamine-transporting, inhibitors; prepn. of polypharmacophoric agents)

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(dopaminergic; prepn. of polypharmacophoric agents)

IT Monoamines

RL: BPR (Biological process); BSU (Biological study, unclassified); BIOL (Biological study); PROC (Process)

(inhibitors; prepn. of polypharmacophoric agents)

IT Transport proteins

RL: BPR (Biological process); BSU (Biological study, unclassified); BIOL (Biological study); PROC (Process)

(monoamine-transporting, inhibitors; prepn. of polypharmacophoric agents)

IT Anti-inflammatory agents  
Antidepressants  
Antiobesity agents  
Pharmacophores  
(prepn. of polypharmacophoric agents)  
IT Alzheimer's disease  
(treatment; prepn. of polypharmacophoric agents)  
IT 9001-66-5, monoamine oxidase 9012-25-3, Catechol O-methyl transferase  
RL: BPR (Biological process); BSU (Biological study, unclassified); BIOL  
(Biological study); PROC  
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IT 67469-69-6P 350583-53-8P 350583-56-1P 350583-58-3P 350583-59-4P  
350583-60-7P 350583-61-8P  
350583-62-9P 350583-63-0P 350583-64-1P 350583-65-2P 350583-66-3P  
350583-67-4P  
RL: BAC (Biological activity or effector, except adverse); BSU (Biological  
study, unclassified); SPN  
(Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP  
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(prepn. of polypharmacophoric agents)

=>